This listing of claims will replace all prior versions and listings of claims in the application.

Listing of claims:

Claim 1 (currently amended): A conformationally restricted polyamine analog of the formula:

E-NH-B-A-B-NH-B-A-B-NH-B-A-B-NH-E

wherein each A is independently selected from the group consisting of: a single bond, C₁-C₆ alkyl, and or C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₃-C₆ cycloalkyl, C₃-C₆ cycloaryl, and C₃-C₆ cycloalkenyl;

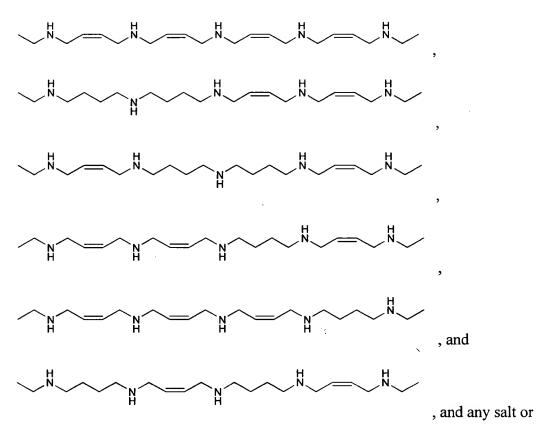
each B is independently selected from the group consisting of: a single bond, C_1 - C_6 alkyl, and C_2 - C_6 alkenyl;

and each E is independently selected from the group consisting of H, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_3 - C_6 cycloalkyl, C_3 - C_6 cycloaryl, and C_3 - C_6 cycloalkenyl;

with the proviso that either at least one A moiety is selected from the group consisting of C_2 - C_6 alkenyl, C_3 - C_6 cycloalkyl, C_3 - C_6 cycloalkyl, and C_3 - C_6 cycloalkenyl, or at least one B moiety is selected from the group consisting of C_2 - C_6 -alkenyl;

and any salt or stereoisomer thereof.

Claim 2 (original): A conformationally restricted polyamine analog according to claim 1, selected from the group consisting of



stereoisomer thereof.

Claim 3 (currently amended): A conformationally restricted polyamine analog of the formula:

 $E-NH-B-A-B-NH-B-A-B-NH(-B-A-B-NH)_x-E$

wherein each A is independently selected from the group consisting of: a single bond, C_1 - C_6 alkyl, and C_2 - C_6 alkenyl, C_3 - C_6 cycloalkyl, C_3 - C_6 cycloalkyl, C_3 - C_6 cycloalkenyl;

each B is independently selected from the group consisting of: a single bond, C_1 - C_6 alkyl, and C_2 - C_6 alkenyl;

each E is independently selected from the group consisting of H, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_3 - C_6 cycloalkyl, C_3 - C_6 cycloaryl, and C_3 - C_6 cycloalkenyl; and x is an integer from 2 to 16;

with the proviso that either at least one A moiety is selected from the group consisting of C_2 - C_6 alkenyl, C_3 - C_6 cycloalkyl, C_3 - C_6 cycloalkyl, and C_3 - C_6 cycloalkenyl, or at least one B moiety is selected from the group consisting of C_2 - C_6 alkenyl;

and any salt or stereoisomer thereof.

Claim 4 (original): A conformationally restricted polyamine analog according to claim 3, selected from the group consisting of:

and

$$\begin{pmatrix} H & H \\ N & N \\ H & N \\ H$$

and any salt or stereoisomer thereof.

Claim 5 (currently amended): A polyamine analog of the formula:

E-NH-B-A-B-NH-B-A-B-NH(-B-A-B-NH)_x-E

wherein each A is independently selected from the group consisting of: a single bond, and C_2 - C_6 alkenyl, C_3 - C_6 -cycloalkyl, C_3 - C_6 -cycloalkyl, C_3 - C_6 -cycloalkyl, and C_3 - C_6 -cycloalkenyl;

each B is independently selected from the group consisting of: a single bond, C_1 - C_6 alkyl, and C_2 - C_6 alkenyl;

each E is independently selected from the group consisting of C₁-C₆ alkyl, C₁-C₆ alkanol, C₃-C₆ cycloalkanol, and C₃-C₆ hydroxyaryl,

with the proviso that at least one E moiety be selected from the group consisting of C_1 - C_6 alkanol, C_3 - C_6 cycloalkanol, and C_3 - C_6 hydroxyaryl;

and x is an integer from 0 to 16;

and any salt or stereoisomer thereof.

Claim 6 (original): A polyamine analog according to claim 5, selected from the group consisting of:

and all salts and stereoisomers thereof.

Claims 7-15 (canceled)

Claim 16 (original): The polyamine analog of claim 1, further comprising a pharmaceutically acceptable excipient.

Claims 17-31 (withdrawn)

Claim 32 (new): The method of claim 17, wherein the indication is breast cancer.